












AMENDMENTS TO THE SPECIFICATION

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









24

Table 1-1 X = Single Bond, q = 0, r = 0, Y = -(R4)C=C(R5)-							
Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R6	R7
1-1-1		H	H	H	H	H	H
1-1-2		H	H	H	H	NO2	H
1-1-3		H	H	H	H	Me	H
1-1-4		H	H	H	H	Cl	H
1-1-5		H	H	H	H	F	H
1-1-6		H	H	H	H	CF3	H
1-1-7		H	H	H	H	COOH	H
1-1-8		H	H	H	Cl	Cl	H
1-1-9		H	H	H	H	H	Me
1-1-10		H	H	H	H	MeO	H
1-1-11		H	H	H	H	H	NO2

Please delete page 37 and replace with the following page 37:























37

(Table 1-2) X = Single Bond, q = 0, r = 0, Y = -(R4)C=C(R5)-

Compound No.	R1-(CH2)r-	R2	R3	R4	R5	R6	R7
1-2-1		H	H	H	H	H	H
1-2-2		H	H	H	H	H	H
1-2-3		H	H	H	H	H	H
1-2-4		H	H	H	H	H	H
1-2-5		H	H	H	H	H	H
1-2-6		H	H	H	H	H	H
1-2-7		H	H	H	H	H	H
1-2-8		H	H	H	H	H	H
1-2-9		H	H	H	H	H	H
1-2-10		H	H	H	H	H	H























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54

(Table 1-3) X = Single Bond, q = 0, r = 0, Y = -(R4)C=O(R5)-							
Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R7	
1-3-1		H	H	H	H	H	
1-3-2		H	H	H	H	H	
1-3-3		H	H	H	H	H	
1-3-4		H	H	H	H	H	
1-3-5		H	H	H	H	H	
1-3-6		H	H	H	H	H	
1-3-7		H	H	H	H		H
1-3-8		H	H	H	H		H
1-3-9		H	H	H	H	H	
1-3-10		H	H	H	H	H	
1-3-11		H	H	H	H	H	












Please delete page 81 and replace with the following page 81:

81

Table 1-4 X = Single Bond, a = 0, r = 0, Y = -(R4)O-(CR5)-							
Compound No.	R1-(CH2)a-	R2	R3	R4	R5	R6	R7
1-4-1		H	H	H	H	H	
1-4-2		H	H	H	H	H	
1-4-3		H	H	H	H	H	
1-4-4		H	H	H	H	H	
1-4-5		H	H	H	H	H	
1-4-6		H	H	H	H	H	
1-4-7		H	H	H	H	H	
1-4-8		H	H	H	H	H	
1-4-9		H	H	H	H	H	
1-4-10		H	H	H	H	H	
1-4-11		H	H	H	H	H	











Please delete page 88 and replace with the following page 88:

88

(Table 1-5) X = Single Bond, q = 0, r = 0, Y = -(R4C-C(R5)-								
Compound No.	R1-(CH2)p-	R2	R3	R4	R5	R6	R7	
1-5-1		H	H	H	H	COOEt	H	
1-5-2		H	H	H	H	COOEt	H	
1-5-3		H	H	H	H	COOEt	H	
1-5-4		H	H	H	H	COOEt	H	
1-5-5		H	H	H	H	COOCH(Me)2	H	
1-5-6		H	H	H	H	COOCH(Me)2	H	
1-5-7		H	H	H	H	COOCH(Me)2	H	
1-5-8		H	H	H	H	COOCH(Me)2	H	
1-5-9		H	H	H	H	COOMe	H	
1-5-10		H	H	H	H	COOMe	H	
1-5-11		H	H	H	H	COOMe	H	

Please delete page 107 and replace with the following page 107:

107

(Table 1-6) X = Single Bond, q = 0, r = 0, Y = -(R4)C=C(R5)-							
Compound No.	R1-(CH2)p--	R2	R3	R4	R5	R6	R7
1-6-1		H	Et	H	H	H	H
1-6-2		Et	Et	H	H	H	H
1-6-3		CH2C6H5	CH2C6H5	H	H	H	H
1-6-4		H	CH2C6H5	H	H	H	H
1-6-5		(CH2)5CH3	(CH2)5CH3	H	H	H	H
1-6-6		H	(CH2)5CH3	H	H	H	H
1-6-7		(CH2)3C6H5	(CH2)3C6H5	H	H	H	H
1-6-8		H	(CH2)3C6H5	H	H	H	H
1-6-9		H	CH2COOMe	H	H	H	H
1-6-10		H	(CH2)4COOEt	H	H	H	H